

# Benzene, 4-chloro-1-(dimethoxymethyl)-

<b>Inchi:</b>	InChI=1S/C9H11ClO2/c1-11-9(12-2)7-3-5-8(10)6-4-7/h3-6,9H,1-2H3
<b>InchiKey:</b>	YRNBYTFFWMWTGG-UHFFFAOYSA-N
<b>Formula:</b>	C9H11ClO2
<b>SMILES:</b>	COC(OC)c1ccc(Cl)cc1
<b>Mol. weight [g/mol]:</b>	186.63
<b>CAS:</b>	3395-81-1

## Physical Properties

Property code	Value	Unit	Source
gf	-96.69	kJ/mol	Joback Method
hf	-289.49	kJ/mol	Joback Method
hfus	15.77	kJ/mol	Joback Method
hvap	47.38	kJ/mol	Joback Method
log10ws	-2.51		Crippen Method
logp	2.631		Crippen Method
mcvol	137.890	ml/mol	McGowan Method
pc	2989.32	kPa	Joback Method
tb	518.81	K	Joback Method
tc	734.53	K	Joback Method
tf	289.51	K	Joback Method
vc	0.510	m3/kmol	Joback Method

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	289.12	J/molxK	518.81	Joback Method
cpg	301.86	J/molxK	554.76	Joback Method
cpg	313.97	J/molxK	590.72	Joback Method
cpg	325.47	J/molxK	626.67	Joback Method
cpg	336.34	J/molxK	662.62	Joback Method
cpg	346.59	J/molxK	698.58	Joback Method
cpg	356.22	J/molxK	734.53	Joback Method
dvisc	0.0018544	Paxs	289.51	Joback Method
dvisc	0.0009761	Paxs	327.73	Joback Method

dvisc	0.0005875	Paxs	365.94	Joback Method
dvisc	0.0003892	Paxs	404.16	Joback Method
dvisc	0.0002769	Paxs	442.38	Joback Method
dvisc	0.0002079	Paxs	480.59	Joback Method
dvisc	0.0001629	Paxs	518.81	Joback Method

## Sources

<b>NIST Webbook:</b>	<a href="http://webbook.nist.gov/cgi/cbook.cgi?ID=C3395811&amp;Units=SI">http://webbook.nist.gov/cgi/cbook.cgi?ID=C3395811&amp;Units=SI</a>
<b>Crippen Method:</b>	<a href="http://pubs.acs.org/doi/abs/10.1021/ci9903071">http://pubs.acs.org/doi/abs/10.1021/ci9903071</a>
<b>Crippen Method:</b>	<a href="https://www.chemeo.com/doc/models/crippen_log10ws">https://www.chemeo.com/doc/models/crippen_log10ws</a>
<b>Joback Method:</b>	<a href="https://en.wikipedia.org/wiki/Joback_method">https://en.wikipedia.org/wiki/Joback_method</a>
<b>McGowan Method:</b>	<a href="http://link.springer.com/article/10.1007/BF02311772">http://link.springer.com/article/10.1007/BF02311772</a>

## Legend

<b>cpg:</b>	Ideal gas heat capacity
<b>dvisc:</b>	Dynamic viscosity
<b>gf:</b>	Standard Gibbs free energy of formation
<b>hf:</b>	Enthalpy of formation at standard conditions
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>hvap:</b>	Enthalpy of vaporization at standard conditions
<b>log10ws:</b>	Log10 of Water solubility in mol/l
<b>logp:</b>	Octanol/Water partition coefficient
<b>mcvol:</b>	McGowan's characteristic volume
<b>pc:</b>	Critical Pressure
<b>tb:</b>	Normal Boiling Point Temperature
<b>tc:</b>	Critical Temperature
<b>tf:</b>	Normal melting (fusion) point
<b>vc:</b>	Critical Volume

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